

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP090324\
 Data File : BP021797.D
 Acq On : 03 Sep 2024 21:22
 Operator : RC/JU
 Sample : P3785-02
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 COSX0

Manual Integrations
 APPROVED

Reviewed By :Jagrut Upadhyay 09/04/2024
 Supervised By :mohammad ahmed 09/07/2024

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 ahmed
 09/07/2024

Quant Time: Sep 04 05:05:48 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP082324.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Aug 29 23:56:59 2024
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.781	152	280967	20.000	ng/u1	0.00
20) Naphthalene-d8	10.563	136	1071696	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.422	164	667371	20.000	ng/u1	-0.01
64) Phenanthrene-d10	17.216	188	1448959	20.000	ng/u1	-0.01
79) Chrysene-d12	21.668	240	1514856	20.000	ng/u1	0.00
88) Perylene-d12	25.062	264	1770194	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.275	96	32176	3.802	ng/uL	0.00
4) Pyridine-d5	3.687	84	102054	4.169	ng/u1	0.00
7) Phenol-d5	6.952	99	255944	8.536	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.110	67	372638	22.836	ng/u1	0.00
11) 2-Chlorophenol-d4	7.316	132	423702	21.887	ng/u1	0.00
15) 4-Methylphenol-d8	8.481	113	339102	14.635	ng/u1	0.00
21) Nitrobenzene-d5	8.922	128	210892	24.459	ng/u1	-0.01
24) 2-Nitrophenol-d4	9.646	143	206514	23.596	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.187	165	365097m	21.169	ng/u1	0.00
31) 4-Chloroaniline-d4	10.698	131	149870	5.973	ng/u1	0.00
46) Dimethylphthalate-d6	13.828	166	1225748	24.003	ng/u1	-0.01
49) Acenaphthylene-d8	14.110	160	1334871	23.283	ng/u1	-0.01
54) 4-Nitrophenol-d4	14.633	143	48063m	4.947	ng/u1	0.00
60) Fluorene-d10	15.428	176	1070993	24.946	ng/u1	-0.02
65) 4,6-Dinitro-2-methylph...	15.545	200	168502	21.144	ng/u1	-0.01
73) Anthracene-d10	17.327	188	1926292	28.012	ng/u1	-0.01
81) Pyrene-d10	19.692	212	2428833	28.048	ng/u1	0.00
92) Benzo(a)pyrene-d12	24.827	264	2654704	28.124	ng/u1	-0.01

Target Compounds Qvalue

(#) = qualifier out of range (m) = manual integration (+) = signals summed

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