

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP082924\
 Data File : BP021732.D
 Acq On : 29 Aug 2024 22:34
 Operator : RC/JU
 Sample : P3721-02
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 DCXY6

Manual Integrations
 APPROVED

Reviewed By :Jagrut Upadhyay 08/30/2024
 Supervised By :mohammad ahmed 09/02/2024

Quant Time: Aug 30 00:08:48 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	7.792	152	476417	20.000	ng/u1	0.00
20) Naphthalene-d8	10.575	136	1978741	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.427	164	1220648	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.233	188	2558631	20.000	ng/u1	0.00
79) Chrysene-d12	21.686	240	2200719	20.000	ng/u1	0.02
88) Perylene-d12	25.092	264	2502038	20.000	ng/u1	0.03
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.275	96	46154	3.216	ng/uL	0.00
4) Pyridine-d5	3.687	84	577076	13.903	ng/u1	0.00
7) Phenol-d5	6.957	99	1203428	23.669	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.122	67	672296	24.297	ng/u1	0.00
11) 2-Chlorophenol-d4	7.322	132	800308	24.381	ng/u1	0.00
15) 4-Methylphenol-d8	8.487	113	770809	19.619	ng/u1	0.00
21) Nitrobenzene-d5	8.934	128	407900	25.622	ng/u1	0.00
24) 2-Nitrophenol-d4	9.657	143	431692	26.714	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.192	165	802662	25.206	ng/u1	0.00
31) 4-Chloroaniline-d4	10.704	131	293889	6.344	ng/u1	0.00
46) Dimethylphthalate-d6	13.827	166	2476580	26.516	ng/u1	-0.01
49) Acenaphthylene-d8	14.116	160	2816540	26.859	ng/u1	0.00
54) 4-Nitrophenol-d4	14.633	143	420221	23.648	ng/u1	0.00
60) Fluorene-d10	15.439	176	2101588	26.764	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.557	200	259579	18.446	ng/u1	0.00
73) Anthracene-d10	17.327	188	3203468	26.381	ng/u1	-0.01
81) Pyrene-d10	19.709	212	3660185	29.095	ng/u1	0.02
92) Benzo(a)pyrene-d12	24.850	264	3644907	27.320	ng/u1	0.01
Target Compounds						
16) Acetophenone	8.610	105	66676m	1.059	ng/u1	Qvalue

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

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