

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092523\
 Data File : BP017339.D
 Acq On : 26 Sep 2023 01:32
 Operator : MA/JU
 Sample : 04472-20
 Misc :
 ALS Vial : 30 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 EZYM7

Manual Integrations
 APPROVED

Reviewed By :Yogesh Patel 09/26/2023
 Supervised By :mohammad ahmed 09/29/2023

Quant Time: Sep 26 02:09:28 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092023.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Sep 21 23:51:07 2023
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.928	152	203238	20.000	ng/u1	-0.01
20) Naphthalene-d8	10.752	136	894785	20.000	ng/u1	-0.01
38) Acenaphthene-d10	14.587	164	569992	20.000	ng/u1	-0.01
64) Phenanthrene-d10	17.357	188	1337202	20.000	ng/u1	-0.01
79) Chrysene-d12	21.463	240	1325014	20.000	ng/u1	-0.01
88) Perylene-d12	23.974	264	1488140	20.000	ng/u1	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.299	96	21714	4.388	ng/uL	0.00
4) Pyridine-d5	3.734	84	292861	21.171	ng/u1	0.00
7) Phenol-d5	7.081	99	230403	13.803	ng/u1	-0.02
9) Bis-(2-Chloroethyl)eth...	7.269	67	243081	24.131	ng/u1	-0.01
11) 2-Chlorophenol-d4	7.446	132	125837	9.505	ng/u1	-0.02
15) 4-Methylphenol-d8	8.640	113	279030	21.502	ng/u1	-0.02
21) Nitrobenzene-d5	9.122	128	154395	23.868	ng/u1	-0.02
24) 2-Nitrophenol-d4	9.840	143	47602	6.777	ng/u1	-0.01
28) 2,4-Dichlorophenol-d3	10.363	165	105735	7.968	ng/u1	-0.02
31) 4-Chloroaniline-d4	10.916	131	472408	24.575	ng/u1	-0.01
46) Dimethylphthalate-d6	14.004	166	1066001	26.107	ng/u1	-0.01
49) Acenaphthylene-d8	14.287	160	1195820	25.463	ng/u1	-0.01
54) 4-Nitrophenol-d4	14.834	143	13311m	1.871	ng/u1	0.00
60) Fluorene-d10	15.581	176	977181	27.798	ng/u1	-0.02
65) 4,6-Dinitro-2-methylph...	15.716	200	7952m	1.055	ng/u1	-0.01
73) Anthracene-d10	17.457	188	1535870	25.897	ng/u1	-0.01
81) Pyrene-d10	19.698	212	1972911	27.534	ng/u1	-0.02
92) Benzo(a)pyrene-d12	23.810	264	1944719	27.142	ng/u1	-0.02

Target Compounds Qvalue

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

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