

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042224\
 Data File : BP020083.D
 Acq On : 22 Apr 2024 19:38
 Operator : MA/JU
 Sample : P2104-20
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0AL7

Quant Time: Apr 23 01:36:42 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP041924.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Apr 23 01:17:52 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.640	152	131690	20.000	ng/u1	0.00
20) Naphthalene-d8	10.416	136	524785	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.322	164	295233	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.169	188	528800	20.000	ng/u1	0.01
79) Chrysene-d12	21.668	240	490413	20.000	ng/u1	0.01
88) Perylene-d12	25.062	264	557722	20.000	ng/u1	0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.228	96	6837	2.127	ng/uL	0.00
4) Pyridine-d5	3.628	84	128	0.014	ng/u1	0.00
7) Phenol-d5	6.863	99	510	0.044	ng/u1	0.04
9) Bis-(2-Chloroethyl)eth...	6.993	67	113744	14.178	ng/u1	0.00
11) 2-Chlorophenol-d4	7.181	132	826	0.105	ng/u1	0.00
15) 4-Methylphenol-d8	8.328	113	6	0.001	ng/u1	-0.02
21) Nitrobenzene-d5	8.804	128	55546	14.715	ng/u1	0.00
24) 2-Nitrophenol-d4	9.510	143	55911	13.702	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.140	165	66	0.008	ng/u1	0.10
31) 4-Chloroaniline-d4	10.581	131	4	0.000	ng/u1	0.01
46) Dimethylphthalate-d6	13.734	166	372314	18.453	ng/u1	0.00
49) Acenaphthylene-d8	14.010	160	2250	0.094	ng/u1	0.01
54) 4-Nitrophenol-d4	14.545	143	880	0.287	ng/u1	-0.01
60) Fluorene-d10	15.351	176	24548	1.367	ng/u1	0.01
65) 4,6-Dinitro-2-methylph...	15.510	200	12963	4.196	ng/u1	0.02
73) Anthracene-d10	17.169	188	528800	22.321	ng/u1	-0.09
81) Pyrene-d10	19.680	212	2678	0.102	ng/u1	0.02
92) Benzo(a)pyrene-d12	24.827	264	2897	0.106	ng/u1	0.02

Target Compounds	Qvalue

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

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