

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP070524\
 Data File : BP020814.D
 Acq On : 06 Jul 2024 10:30
 Operator : MA/JU
 Sample : PB161717BL
 Misc :
 ALS Vial : 34 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK717

Quant Time: Jul 07 22:55:49 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062524.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Jul 03 15:01:56 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.728	152	238781	20.000	ng/u1	0.00
20) Naphthalene-d8	10.499	136	999894	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.357	164	663022	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.169	188	1429363	20.000	ng/u1	0.00
79) Chrysene-d12	21.628	240	1379803	20.000	ng/u1	0.00
88) Perylene-d12	24.974	264	1583982	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.264	96	27561	4.956	ng/uL	0.00
4) Pyridine-d5	3.676	84	400982	24.743	ng/u1	0.00
7) Phenol-d5	6.917	99	522131	26.660	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.069	67	303135	24.415	ng/u1	-0.01
11) 2-Chlorophenol-d4	7.275	132	424545	26.524	ng/u1	0.00
15) 4-Methylphenol-d8	8.440	113	420896	26.461	ng/u1	0.00
21) Nitrobenzene-d5	8.881	128	212402	28.772	ng/u1	0.00
24) 2-Nitrophenol-d4	9.587	143	243286	32.932	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.134	165	420975	27.664	ng/u1	0.00
31) 4-Chloroaniline-d4	10.640	131	579748	27.185	ng/u1	0.00
46) Dimethylphthalate-d6	13.769	166	1321170	28.659	ng/u1	0.01
49) Acenaphthylene-d8	14.046	160	1475377	26.821	ng/u1	0.00
54) 4-Nitrophenol-d4	14.599	143	211479	30.025	ng/u1	0.02
60) Fluorene-d10	15.363	176	1115574	28.758	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.493	200	157837	21.912	ng/u1	0.00
73) Anthracene-d10	17.269	188	1830801	28.539	ng/u1	0.00
81) Pyrene-d10	19.657	212	2187300	26.387	ng/u1	0.00
92) Benzo(a)pyrene-d12	24.739	264	2235205	29.009	ng/u1	0.00

Target Compounds Qvalue

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

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