

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072924\
 Data File : BP021236.D
 Acq On : 30 Jul 2024 15:36
 Operator : MA/JU
 Sample : PB162276BL
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK276

Quant Time: Jul 30 22:54:01 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Jul 29 12:46:18 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.628	152	205585	20.000	ng/u1	0.00
20) Naphthalene-d8	10.393	136	809096	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.275	164	543498	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.098	188	1203574	20.000	ng/u1	0.00
79) Chrysene-d12	21.551	240	1166855	20.000	ng/u1	# 0.02
88) Perylene-d12	24.851	264	1217494	20.000	ng/u1	0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.170	96	22326	4.942	ng/uL	0.00
4) Pyridine-d5	3.593	84	284325	21.739	ng/u1	0.00
7) Phenol-d5	6.846	99	368590	21.991	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	6.981	67	220454	21.704	ng/u1	0.00
11) 2-Chlorophenol-d4	7.181	132	321024	23.247	ng/u1	0.00
15) 4-Methylphenol-d8	8.358	113	312607	22.456	ng/u1	0.00
21) Nitrobenzene-d5	8.793	128	156273	24.867	ng/u1	0.00
24) 2-Nitrophenol-d4	9.499	143	179521	24.958	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.052	165	290474	22.333	ng/u1	0.00
31) 4-Chloroaniline-d4	10.552	131	415925	24.139	ng/u1	0.00
46) Dimethylphthalate-d6	13.693	166	1035472	26.208	ng/u1	0.01
49) Acenaphthylene-d8	13.969	160	1114456	25.121	ng/u1	0.00
54) 4-Nitrophenol-d4	14.593	143	81851	14.561	ng/u1	0.00
60) Fluorene-d10	15.293	176	885969	27.334	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.475	200	136502	18.744	ng/u1	0.01
73) Anthracene-d10	17.204	188	1392642	26.052	ng/u1	0.01
81) Pyrene-d10	19.586	212	1773074	24.578	ng/u1	0.01
92) Benzo(a)pyrene-d12	24.621	264	1627319	27.101	ng/u1	0.02

Target Compounds Qvalue

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

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