

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP080324\
 Data File : BP021338.D
 Acq On : 03 Aug 2024 10:13
 Operator : CG/JU
 Sample : PB162169BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK169

Quant Time: Aug 04 23:20:35 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	216097	20.000	ng/u1	0.00
20) Naphthalene-d8	10.381	136	819016	20.000	ng/u1	-0.01
38) Acenaphthene-d10	14.263	164	468268	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.081	188	964082	20.000	ng/u1	-0.01
79) Chrysene-d12	21.522	240	1094738	20.000	ng/u1	-0.01
88) Perylene-d12	24.821	264	1380193	20.000	ng/u1	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.164	96	24246	5.106	ng/uL	-0.01
4) Pyridine-d5	3.587	84	303775	22.096	ng/u1	0.00
7) Phenol-d5	6.852	99	379358	21.533	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	6.975	67	243468	22.803	ng/u1	0.00
11) 2-Chlorophenol-d4	7.175	132	331249	22.820	ng/u1	0.00
15) 4-Methylphenol-d8	8.363	113	312047	21.326	ng/u1	0.00
21) Nitrobenzene-d5	8.787	128	148631	23.365	ng/u1	0.00
24) 2-Nitrophenol-d4	9.505	143	169356	23.260	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.052	165	277852	21.104	ng/u1	0.00
31) 4-Chloroaniline-d4	10.563	131	390663	22.398	ng/u1	0.00
46) Dimethylphthalate-d6	13.675	166	793880	23.322	ng/u1	0.00
49) Acenaphthylene-d8	13.951	160	909430	23.793	ng/u1	-0.01
54) 4-Nitrophenol-d4	14.622	143	72845	15.040	ng/u1	0.04
60) Fluorene-d10	15.281	176	675958	24.205	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.457	200	58359	10.005	ng/u1	0.00
73) Anthracene-d10	17.187	188	1101170	25.716	ng/u1	0.00
81) Pyrene-d10	19.569	212	1410420	20.839	ng/u1	0.00
92) Benzo(a)pyrene-d12	24.592	264	1722106	25.299	ng/u1	0.00

Target Compounds Qvalue

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

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