

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP070323\
 Data File : BP015959.D
 Acq On : 03 Jul 2023 20:01
 Operator : MA/JU
 Sample : PB153768BL
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK768

Quant Time: Jul 04 00:00:35 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062623.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Sun Jul 02 06:12:16 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.046	152	303572	20.000	ng/u1	0.00
20) Naphthalene-d8	10.875	136	1359456	20.000	ng/u1	# 0.00
38) Acenaphthene-d10	14.698	164	814393	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.457	188	1650103	20.000	ng/u1	# 0.00
79) Chrysene-d12	21.557	240	1102404	20.000	ng/u1	# 0.00
88) Perylene-d12	24.098	264	1169946	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.370	96	45944	5.445	ng/uL	0.00
4) Pyridine-d5	3.811	84	631925	29.711	ng/u1	0.00
7) Phenol-d5	7.228	99	820821	31.602	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.369	67	566464	35.251	ng/u1	0.00
11) 2-Chlorophenol-d4	7.575	132	659105	33.616	ng/u1	0.00
15) 4-Methylphenol-d8	8.787	113	692248	33.737	ng/u1	0.00
21) Nitrobenzene-d5	9.234	128	335667	32.308	ng/u1	0.00
24) 2-Nitrophenol-d4	9.963	143	367637	30.319	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.534	165	595183	27.544	ng/u1	0.02
31) 4-Chloroaniline-d4	11.040	131	874422	31.502	ng/u1	0.01
46) Dimethylphthalate-d6	14.110	166	2009875	31.930	ng/u1	0.00
49) Acenaphthylene-d8	14.393	160	2259573	31.933	ng/u1	0.00
54) 4-Nitrophenol-d4	14.998	143	217265	26.156	ng/u1	0.05
60) Fluorene-d10	15.698	176	1652099	31.875	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.869	200	201099	19.817	ng/u1	0.01
73) Anthracene-d10	17.563	188	2481928	32.928	ng/u1	0.00
81) Pyrene-d10	19.792	212	2533747	35.198	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.927	264	1983244	33.604	ng/u1	0.00

Target Compounds Qvalue

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

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