

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP071922\
 Data File : BP011031.D
 Acq On : 19 Jul 2022 14:23
 Operator : CG/JU
 Sample : PB146296BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK296

Quant Time: Jul 19 23:40:19 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071422.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Jul 19 03:25:54 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.675	152	356025	20.000	ng/u1	0.00
20) Naphthalene-d8	10.469	136	1435819	20.000	ng/u1	# 0.00
38) Acenaphthene-d10	14.340	164	799162	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.110	188	1630487	20.000	ng/u1	# 0.00
79) Chrysene-d12	21.228	240	1571161	20.000	ng/u1	# 0.00
88) Perylene-d12	23.586	264	1581351	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.181	96	62151	6.293	ng/uL	0.00
4) Pyridine-d5	3.587	84	773438	30.300	ng/u1	0.00
7) Phenol-d5	6.858	99	876056	29.976	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.022	67	586880	31.229	ng/u1	0.00
11) 2-Chlorophenol-d4	7.211	132	743530	31.486	ng/u1	0.00
15) 4-Methylphenol-d8	8.393	113	698756	29.866	ng/u1	0.00
21) Nitrobenzene-d5	8.840	128	337921	32.295	ng/u1	0.00
24) 2-Nitrophenol-d4	9.558	143	333812	32.241	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.093	165	576743	30.460	ng/u1	0.00
31) 4-Chloroaniline-d4	10.616	131	959491	30.738	ng/u1	0.00
46) Dimethylphthalate-d6	13.757	166	1856743	33.994	ng/u1	0.00
49) Acenaphthylene-d8	14.028	160	2243087	34.806	ng/u1	0.00
54) 4-Nitrophenol-d4	14.557	143	278089	25.156	ng/u1	0.00
60) Fluorene-d10	15.340	176	1560514	33.612	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.463	200	189539	23.492	ng/u1	0.00
73) Anthracene-d10	17.204	188	2364062	34.106	ng/u1	0.00
81) Pyrene-d10	19.463	212	2686972	32.562	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.433	264	2563584	33.753	ng/u1	0.00

Target Compounds Qvalue

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

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