

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP080522\
 Data File : BP011273.D
 Acq On : 05 Aug 2022 11:24
 Operator : CG/JU
 Sample : PB146741BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK741

Quant Time: Aug 05 23:01:07 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP080222.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Aug 02 23:52:37 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.599	152	120806	20.000	ng/u1	0.00
20) Naphthalene-d8	10.393	136	507480	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.275	164	269116	20.000	ng/u1	# 0.00
64) Phenanthrene-d10	17.045	188	516538	20.000	ng/u1	0.00
79) Chrysene-d12	21.169	240	646930	20.000	ng/u1	0.00
88) Perylene-d12	23.486	264	703855	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.087	96	6124	4.099	ng/uL	0.00
4) Pyridine-d5	3.505	84	69349	17.208	ng/u1	0.00
7) Phenol-d5	6.781	99	296093	29.411	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	6.946	67	279707	38.503	ng/u1	0.00
11) 2-Chlorophenol-d4	7.134	132	246802	30.076	ng/u1	0.00
15) 4-Methylphenol-d8	8.322	113	233956	28.243	ng/u1	0.00
21) Nitrobenzene-d5	8.763	128	164792	45.146	ng/u1	0.00
24) 2-Nitrophenol-d4	9.481	143	114263	33.395	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.022	165	205340	31.472	ng/u1	0.00
31) 4-Chloroaniline-d4	10.546	131	288053	28.537	ng/u1	0.00
46) Dimethylphthalate-d6	13.698	166	652693	35.411	ng/u1	0.00
49) Acenaphthylene-d8	13.963	160	1015234	43.778	ng/u1	0.00
54) 4-Nitrophenol-d4	14.504	143	77253	24.648	ng/u1	0.00
60) Fluorene-d10	15.275	176	776803	49.530	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.410	200	72851	30.374	ng/u1	0.00
73) Anthracene-d10	17.139	188	1365428	60.147	ng/u1	0.00
81) Pyrene-d10	19.404	212	1473706	45.055	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.339	264	1553189	44.283	ng/u1	0.00

Target Compounds Qvalue

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

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