

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041322\
 Data File : BP009812.D
 Acq On : 13 Apr 2022 11:41
 Operator : CG/JU
 Sample : PB144076BL
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK076

Quant Time: Apr 14 01:05:03 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP040722.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Apr 13 02:15:22 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.963	152	150575	20.000	ng/u1	0.00
20) Naphthalene-d8	10.769	136	684233	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.593	164	490684	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.345	188	1091414	20.000	ng/u1	# 0.00
79) Chrysene-d12	21.422	240	1174355	20.000	ng/u1	#-0.01
88) Perylene-d12	23.892	264	1098023	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.387	96	20900	6.195	ng/uL	0.00
4) Pyridine-d5	3.799	84	281341	30.244	ng/u1	0.00
7) Phenol-d5	7.116	99	363626	27.507	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.287	67	247474	31.445	ng/u1	0.00
11) 2-Chlorophenol-d4	7.493	132	297225	29.807	ng/u1	0.00
15) 4-Methylphenol-d8	8.663	113	318671	28.877	ng/u1	0.00
21) Nitrobenzene-d5	9.122	128	159048	32.234	ng/u1	0.00
24) 2-Nitrophenol-d4	9.846	143	175259	32.398	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.387	165	310848	27.460	ng/u1	0.00
31) 4-Chloroaniline-d4	10.899	131	469584	29.125	ng/u1	0.00
46) Dimethylphthalate-d6	14.004	166	1176994	30.536	ng/u1	0.00
49) Acenaphthylene-d8	14.287	160	1292109	28.296	ng/u1	0.00
54) 4-Nitrophenol-d4	14.775	143	202558	28.978	ng/u1	0.00
60) Fluorene-d10	15.587	176	989257	30.510	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.692	200	174884	27.049	ng/u1	0.00
73) Anthracene-d10	17.445	188	1598409	30.588	ng/u1	0.00
81) Pyrene-d10	19.675	212	1984327	30.939	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.733	264	1824512	31.908	ng/u1	0.00

Target Compounds Qvalue

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

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