

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041922\
 Data File : BP009995.D
 Acq On : 21 Apr 2022 02:15
 Operator : CG/JU
 Sample : PB144244BL
 Misc :
 ALS Vial : 49 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK244

Quant Time: Apr 21 03:03:34 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP040722.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Apr 20 01:32:14 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.946	152	112322	20.000	ng/u1	0.00
20) Naphthalene-d8	10.746	136	515614	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.575	164	375941	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.322	188	837313	20.000	ng/u1	# 0.00
79) Chrysene-d12	21.404	240	878076	20.000	ng/u1	0.00
88) Perylene-d12	23.851	264	826210	20.000	ng/u1	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.370	96	17714	7.038	ng/uL	0.00
4) Pyridine-d5	3.781	84	216991	31.271	ng/u1	0.00
7) Phenol-d5	7.099	99	304797	30.909	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.269	67	210018	35.774	ng/u1	0.00
11) 2-Chlorophenol-d4	7.475	132	253862	34.128	ng/u1	0.00
15) 4-Methylphenol-d8	8.646	113	268790	32.652	ng/u1	0.00
21) Nitrobenzene-d5	9.099	128	137544	36.992	ng/u1	0.00
24) 2-Nitrophenol-d4	9.828	143	150116	36.825	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.363	165	258416	30.294	ng/u1	0.00
31) 4-Chloroaniline-d4	10.881	131	383581	31.571	ng/u1	0.00
46) Dimethylphthalate-d6	13.981	166	1044136	35.358	ng/u1	0.00
49) Acenaphthylene-d8	14.263	160	1133524	32.400	ng/u1	0.00
54) 4-Nitrophenol-d4	14.757	143	157936	29.490	ng/u1	0.00
60) Fluorene-d10	15.563	176	869802	35.013	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.681	200	143577	28.945	ng/u1	0.00
73) Anthracene-d10	17.422	188	1414651	35.287	ng/u1	0.00
81) Pyrene-d10	19.651	212	1740606	36.296	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.698	264	1526062	35.468	ng/u1	-0.01

Target Compounds Qvalue

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

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