

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041122\
 Data File : BP009787.D
 Acq On : 12 Apr 2022 17:34
 Operator : CG/JU
 Sample : N2316-18
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 ETNR6

Quant Time: Apr 13 02:23:53 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	7.969	152	109599	20.000	ng/u1	0.00
20) Naphthalene-d8	10.775	136	493698	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.598	164	373019	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.345	188	839314	20.000	ng/u1	# 0.00
79) Chrysene-d12	21.427	240	888557	20.000	ng/u1	0.00
88) Perylene-d12	23.898	264	836440	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.387	96	9833	4.004	ng/uL	0.00
4) Pyridine-d5	3.811	84	31603	4.668	ng/u1	0.00
7) Phenol-d5	7.116	99	53988	5.611	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.293	67	170553	29.773	ng/u1	0.00
11) 2-Chlorophenol-d4	7.493	132	159103	21.921	ng/u1	0.00
15) 4-Methylphenol-d8	8.663	113	109026	13.573	ng/u1	0.00
21) Nitrobenzene-d5	9.122	128	111486	31.315	ng/u1	0.00
24) 2-Nitrophenol-d4	9.852	143	110799	28.386	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.387	165	196895	24.106	ng/u1	0.00
31) 4-Chloroaniline-d4	10.904	131	287475	24.712	ng/u1	0.00
46) Dimethylphthalate-d6	14.004	166	886456	30.253	ng/u1	0.00
49) Acenaphthylene-d8	14.292	160	953169	27.458	ng/u1	0.00
54) 4-Nitrophenol-d4	14.769	143	27943	5.258	ng/u1	-0.01
60) Fluorene-d10	15.592	176	764904	31.032	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.698	200	126525	25.447	ng/u1	0.00
73) Anthracene-d10	17.445	188	1382021	34.391	ng/u1	0.00
81) Pyrene-d10	19.675	212	1696497	34.959	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.739	264	1534018	35.217	ng/u1	0.00

Target Compounds	Qvalue

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

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